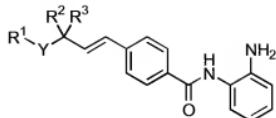


1. - 11. (Canceled)

12. (Previously Presented) A compound of formula:



or a pharmaceutically acceptable salt thereof, wherein

Y is -N(R⁴)-, -O-, -S-, -N(R⁴)SO₂-, -SO₂-N(R⁴)-, -SO₂-, -N(R⁴)-C(O)-, -C(O)-N(R⁴)-, -NHC(O)NH-, -N(R⁴)C(O)O-, -OC(O)N(R⁴)-, or a covalent bond, and

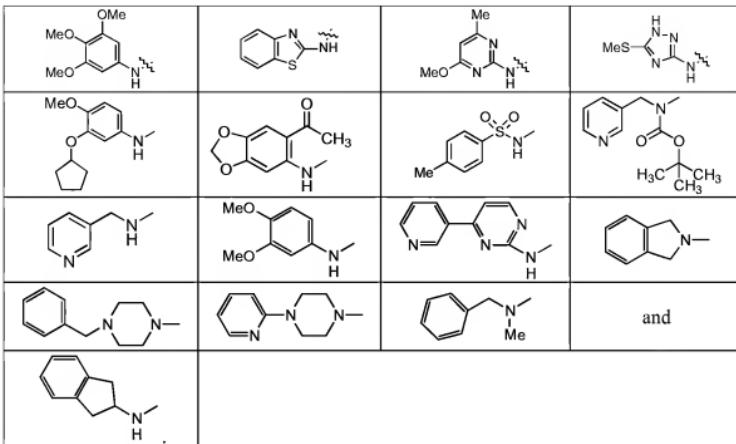
R¹, R², and R³ independently are -H or R^a-C₀-C₆-hydrocarbyl wherein R^a is -H or R^a is aryl or heteroaryl, each of which is optionally substituted with from 1 to 3 substituents.

R⁴ is -H, -C(O)-R^b, -C(O)O-R^b, -C(O)NH-R^b, or R^c-C₀-C₆-hydrocarbyl wherein

R^b is -H or -C₁-C₆-hydrocarbyl, and

R^c is -H, or aryl or heteroaryl each of which is optionally substituted with from 1 to 3 substituents.

13. (Original) The compound according to claim 12 wherein R² and R³ are both -H.
14. (Original) The compound according to claim 12 wherein Y is -NH-, -SO₂-NH-, or -N(R⁴)- wherein R⁴ is -C(O)O-C₁-C₆-hydrocarbyl.
15. (Original) The compound according to claim 12 wherein R¹ is aryl, benzothiazolyl, pyrimidinyl, triazolyl, benzodioxolenyl, or pyridinyl, each of which is optionally substituted with from 1 to 3 substituents.
16. (Original) The compound according to claim 15 wherein R¹ is substituted with from 1-3 substituents independently selected from C₁-C₆-hydrocarbyl, C₁-C₆-hydrocarbyloxy, halo, methylthio, and acetyl.
17. (Previously Presented) The compound according to claim 12 wherein R¹-Y is selected from :



18. – 25. (Canceled)